Welcome to STN International! Enter x:X

LOGINID:ssptajqm1797

PASSWORD: TERMINAL (ENTER 1, 2, 3, OR ?):2

TERMIN	IAL	(ENT	ER 1	, 2, 3, OR ?):2
* * *	* *	* *	* *	* Welcome to STN International * * * * * * * * * *
NEWS	1			Web Page for STN Seminar Schedule - N. America
NEWS	2	NOV	21	CAS patent coverage to include exemplified prophetic substances identified in English-, French-, German-,
	_			and Japanese-language basic patents from 2004-present
NEWS	3	NOV		MARPAT enhanced with FSORT command
NEWS	4 5	NOV		CHEMSAFE now available on STN Easy
	5			Two new SET commands increase convenience of STN searching
NEWS	6	DEC		ChemPort single article sales feature unavailable
NEWS	7	DEC	12	GBFULL now offers single source for full-text coverage of complete UK patent families
NEWS	8	DEC	17	Fifty-one pharmaceutical ingredients added to PS
NEWS	9	JAN	06	The retention policy for unread STNmail messages will change in 2009 for STN-Columbus and STN-Tokyo
NEWS	10	JAN	07	WPIDS, WPINDEX, and WPIX enhanced Japanese Patent
NEWS	11	FEB	02	Simultaneous left and right truncation (SLART) added
NIENTO	10		0.0	for CERAB, COMPUAB, ELCOM, and SOLIDSTATE
NEWS		FEB		GENBANK enhanced with SET PLURALS and SET SPELLING
NEWS		FEB		Patent sequence location (PSL) data added to USGENE
NEWS		FEB		COMPENDEX reloaded and enhanced
NEWS		FEB		WTEXTILES reloaded and enhanced
NEWS	16	FEB	19	New patent-examiner citations in 300,000 CA/CAplus patent records provide insights into related prior art
NEWS	17	FEB	19	Increase the precision of your patent queries use terms from the IPC Thesaurus, Version 2009.01
NEWS	18	FEB	23	Several formats for image display and print options discontinued in USPATFULL and USPAT2
NEWS	19	FEB	23	MEDLINE now offers more precise author group fields and 2009 MeSH terms
NEWS	20	FEB	23	TOXCENTER updates mirror those of MEDLINE - more
NEWS	21	FEB	23	precise author group fields and 2009 MeSH terms Three million new patent records blast AEROSPACE into
NEWS	22	FEB	25	STN patent clusters USGENE enhanced with patent family and legal status
NEWS	23	MAR	06	display data from INPADOCDB INPADOCDB and INPAFAMDB enhanced with new display
NEWS	24	MAR	11	formats EPFULL backfile enhanced with additional full-text
				applications and grants
NEWS		MAR		ESBIOBASE reloaded and enhanced
NEWS	26	MAR	20	CAS databases on STN enhanced with new super role for nanomaterial substances
NEWS	27	MAR	23	${ m CA/CAplus}$ enhanced with more than 250,000 patent equivalents from China

NEWS EXPRESS JUNE 27 08 CURRENT WINDOWS VERSION IS V8.3, AND CURRENT DISCOVER FILE IS DATED 23 JUNE 2008.

NEWS HOURS STN Operating Hours Plus Help Desk Availability

NEWS LOGIN Welcome Banner and News Items

NEWS IPC8 For general information regarding STN implementation of IPC 8

Enter NEWS followed by the item number or name to see news on that specific topic.

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FILE 'HOME' ENTERED AT 11:04:41 ON 24 MAR 2009

=> file caslink

COST IN U.S. DOLLARS

SINCE FILE 0.22

TOTAL ENTRY SESSION 0.22

FULL ESTIMATED COST

FILE 'CAPLUS' ENTERED AT 11:04:57 ON 24 MAR 2009 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

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FILE 'MARPAT' ENTERED AT 11:04:57 ON 24 MAR 2009 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2009 American Chemical Society (ACS)

FILE 'REGISTRY' ENTERED AT 11:04:57 ON 24 MAR 2009 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2009 American Chemical Society (ACS)

CLUSTER 'CASLINK' ENTERED

Predefined command sequences will be executed in REGISTRY, MARPAT, and CAPLUS.

Uploading C:\Program Files\STNEXP\Oueries\10529897-b.str

```
chain nodes: 1 2 3 5 6 7 chain bonds: 1-2 1-3 1-5 1-6 6-7 exact/norm bonds: 1-2 1-3 1-5 exact bonds: 1-6 6-7
```

G2:H,CH3,Et,n-Pr,i-Pr,n-Bu,i-Bu,s-Bu,t-Bu

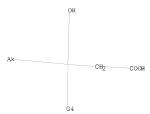
G3:CH3,Et,n-Pr,n-Bu

G4:H,Ak

Match level : 1:CLASS 2:CLASS 3:CLASS 5:CLASS 6:CLASS 7:CLASS

L1 STRUCTURE UPLOADED

=> d L1 HAS NO ANSWERS L1 STR



G1

- G2 H, Me, Et, n-Pr, i-Pr, n-Bu, i-Bu, s-Bu, t-Bu
- G3 Me, Et, n-Pr, n-Bu

G4 H, Ak

Structure attributes must be viewed using STN Express query preparation.

=> s 11 sss sample

S L1 SSS SAM FILE=REGISTRY SAMPLE SEARCH INITIATED 11:05:30 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 45483 TO ITERATE

4.4% PROCESSED 2000 ITERATIONS INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 896919 TO 922401

PROJECTED ITERATIONS: 896919 TO 92 PROJECTED ANSWERS: 22023 TO 2

L2 50 SEA SSS SAM L1 1 FILES SEARCHED...

S L2 SSS SAM FILE-MARPAT SAMPLE SEARCH INITIATED 11:05:31 FILE 'MARPAT' SAMPLE SCREEN SEARCH COMPLETED - 11176 TO ITERATE

17.9% PROCESSED 2000 ITERATIONS INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.02

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
PROJECTED ITERATIONS: 220004 TO 227036
PROJECTED ANSWERS: 16025 TO 19513

L3 50 SEA SSS SAM L1 1 FILES SEARCHED...

=> s 11 sss ful1

50 ANSWERS

50 ANSWERS

S L1 SSS FUL FILE=REGISTRY

FULL SEARCH INITIATED 11:05:50 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 906193 TO ITERATE

90.7% PROCESSED 821687 ITERATIONS

24895 ANSWERS

99.0% PROCESSED 896999 ITERATIONS

25541 ANSWERS

100.0% PROCESSED 906193 ITERATIONS

25559 ANSWERS

SEARCH TIME: 00.00.38

25559 SEA SSS FUL L1 1 FILES SEARCHED...

S L4 SSS FUL FILE=MARPAT

FULL SEARCH INITIATED 11:06:29 FILE 'MARPAT'

FULL SCREEN SEARCH COMPLETED - 223897 TO ITERATE

32.7% PROCESSED 73128 ITERATIONS 7047 ANSWERS

62.2% PROCESSED 139324 ITERATIONS

13469 ANSWERS

77.3% PROCESSED 172965 ITERATIONS 16342 ANSWERS

89.5% PROCESSED 200483 ITERATIONS (17914 ANSWERS

18910 ANSWERS

91.1% PROCESSED 203862 ITERATIONS (3 INCOMPLETE) 18258 ANSWERS

217743 ITERATIONS (97.3% PROCESSED 98.2% PROCESSED 219909 ITERATIONS (

7 INCOMPLETE) 19017 ANSWERS

98.3% PROCESSED 220177 ITERATIONS (98.4% PROCESSED 220367 ITERATIONS (

7 INCOMPLETE) 19056 ANSWERS

8 INCOMPLETE) 19087 ANSWERS 19099 ANSWERS

8 INCOMPLETE) 9 INCOMPLETE)

2 INCOMPLETE)

5 INCOMPLETE)

19147 ANSWERS 19331 ANSWERS

100.0% PROCESSED 223897 ITERATIONS (SEARCH TIME: 00.03.33

99.5% PROCESSED 222706 ITERATIONS (

220499 ITERATIONS (

220862 ITERATIONS (

10 INCOMPLETE) 10 INCOMPLETE) 19455 ANSWERS

19455 SEA SSS FUL L1 1 FILES SEARCHED...

98.5% PROCESSED

98.6% PROCESSED

S L4 FILE=CAPLUS L6 162344 FILE CAPLUS 1 FILES SEARCHED...

SET DUPORDER FILE

SET COMMAND COMPLETED

DUP REM L5 L6

181799 ANSWERS REQUESTED EXCEEDS MAXIMUM ALLOWED OF 50000 You may process up to 50,000 answers per command. Please try to narrow your search until your resulting L# answer set is within the maximum number of answers.

```
=> delete history yes
ALL L# ITEMS DELETED
```

=> delete select yes
ALL E# DEFINITIONS DELETED

=>

Uploading C:\Program Files\STNEXP\Queries\10529897-b-narrow.str



```
chain nodes: 1 2 3 5 6 7 11 chain bonds: 1-2 1-3 1-5 1-6 6-7 exact/norm bonds: 1-2 1-3 1-5 exact bonds: 1-6 6-7
```

G2:H,CH3,Et,n-Pr,i-Pr,n-Bu,i-Bu,s-Bu,t-Bu

G3:CH3,Et,n-Pr,n-Bu

G4:H,[*1]

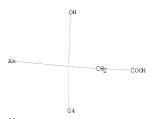
Match level:
1:CLASS 2:CLASS 3:CLASS 5:CLASS 6:CLASS 7:CLASS 11:CLASS
Element Count:
Node 2: Limited
C.Cl-4

Node 11: Limited C,C1-4 I.1 STRUCTURE UPLOADED

=> d

L1 HAS NO ANSWERS

L1 STR



G2 H, Me, Et, n-Pr, i-Pr, n-Bu, i-Bu, s-Bu, t-Bu

G4 H, [@1]

=> s 11 sss sample

S L1 SSS SAM FILE=REGISTRY

SAMPLE SEARCH INITIATED 11:12:49 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 45483 TO ITERATE

4.4% PROCESSED 2000 ITERATIONS INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE** BATCH **COMPLETE**

PROJECTED ITERATIONS: 896919 TO 922401 PROJECTED ANSWERS: 19849 TO 23813

48 SEA SSS SAM L1

1 FILES SEARCHED...

S L2 SSS SAM FILE=MARPAT SAMPLE SEARCH INITIATED 11:12:49 FILE 'MARPAT'

SAMPLE SCREEN SEARCH COMPLETED - 11176 TO ITERATE

17.9% PROCESSED 2000 ITERATIONS INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 220004 TO 227036 Ak 1

48 ANSWERS

50 ANSWERS

G3 Me, Et, n-Pr, n-Bu

Structure attributes must be viewed using STN Express query preparation.

PROJECTED ANSWERS: 16025 TO 19513

L3 50 SEA SSS SAM L1 1 FILES SEARCHED...

=> d scan

L3 50 ANSWERS MARPAT COPYRIGHT 2009 ACS on STN

IC ICM A61K031-663

ICS A61K031-5025; A61K031-662; A61K045-00; A61P001-02; A61P019-00; A61P019-02; A61P019-04; A61P019-08; A61P019-10; A61P029-00; A61P035-00: A61P035-00.

CC 63-6 (Pharmaceuticals)

Section cross-reference(s): 1

TI Agent inducing increase in bone mass

ST bisphosphonate deriv osteoblast differentiation promoter combination bone mass increase; pyridazine deriv prepn incadronate combination bone mass increase

IT Bone

Bone resorption inhibitors

Human

Osteoporosis

(bone mass-increasing agents containing nonpeptidic osteoblast differentiation promoting compds. and bisphosphonates)

IT Drug delivery systems

(capsules; bone mass-increasing agents containing nonpeptidic osteoblast differentiation promoting compds. and bisphosphonates)

IT Osteoblast

(differentiation; bone mass-increasing agents containing nonpeptidic

osteoblast differentiation promoting compds. and bisphosphonates)

IT Cell differentiation

(osteoblast; bone mass-increasing agents containing nonpeptidic osteoblast differentiation promoting compds. and bisphosphonates)

IT Bone, disease

(osteopenia; bone mass-increasing agents containing nonpeptidic osteoblast differentiation promoting compds. and bisphosphonates)

IT Drug delivery systems

(tablets; bone mass-increasing agents containing nonpeptidic osteoblast

differentiation promoting compds. and bisphosphonates)

	differentia	ation promoting	compas. and bi	spnospnonates)	
ΙT	596823-90-4P	596823-92-6P	596823-94-8P	596823-96-0P	596823-98-2P
	596824-04-3P	596824-07-6P	596824-09-8P	596824-12-3P	596824-13-4P
	596824-15-6P	596824-16-7P	596824-17-8P	596824-18-9P	596824-19-0P
	596824-20-3P	596824-21-4P	596824-22-5P	596824-23-6P	596824-24-7P
	596824-25-8P	596824-26-9P	596824-27-0P	596824-28-1P	596824-29-2P
	596824-30-5P	596824-31-6P	596824-32-7P	596824-33-8P	596824-34-9P
	596824-36-1P	596824-38-3P	596824-40-7P	596824-42-9P	596824-43-0P
	596824-44-1P	596824-45-2P	596824-46-3P	596824-47-4P	596824-48-5P
	596824-49-6P	596824-50-9P	596824-51-0P	596824-52-1P	596824-53-2P
	596824-54-3P	596824-55-4P	596824-56-5P	596824-57-6P	596824-58-7P
	596824-59-8P	596824-60-1P	596824-61-2P	596824-62-3P	596824-63-4P
	596824-64-5P	596824-65-6P	596824-66-7P	596824-67-8P	596824-68-9P
	596824-69-0P	596824-70-3P	596824-71-4P	596824-72-5P	596824-73-6P
	596824-74-7P	596824-75-8P	596824-76-9P	596824-77-0P	596824-79-2P
	596824-81-6P	596824-83-8P	596824-84-9P	596824-85-0P	596824-86-1P
	596824-87-2P	596824-88-3P	596824-89-4P	596824-90-7P	596824-91-8P
	596824-92-9P	596824-93-0P	596824-95-2P	596824-96-3P	596824-98-5P
	596825-00-2P	596825-02-4P	596825-03-5P	596825-04-6P	596825-05-7P
	596825-06-8P	596825-07-9P	596825-08-0P	596825-09-1P	596825-10-4P
	596825-11-5P	596825-12-6P	596825-13-7P	596825-14-8P	596825-15-9P
	596825-16-0P	596825-17-1P	596825-18-2P	596825-19-3P	596825-20-6P
	596825-21-7P	596825-22-8P	596825-23-9P	596825-24-0P	596825-25-1P
	596825-26-2P	596825-27-3P	596825-28-4P	596825-29-5P	596825-30-8P

```
596825-31-9P 596825-32-0P 596825-33-1P 596825-34-2P 596825-35-3P RI: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
```

(bone mass-increasing agents containing nonpeptidic osteoblast differentiation promoting compds. and bisphosphonates)

RL: THU (Therapeutic use); BIOL (Biological study); USES (Uses) (bone mass-increasing agents containing nonpeptidic osteoblast differentiation promoting compds. and bisphosphonates)

IT 62-23-7, 4-Mitrobenzoic acid 93-64-9, 3-Dimethylaminobenzoic acid
110-89-4, Piperidine, reactions 111-49-9 124-63-0, Methanesulfonyl
chloride 1121-92-2 1711-11-1, 3-Cyanobenzoylchloride 4684-94-0,
6-Chloropicolinic acid 5398-36-7, Ethyl 2-aminothiazole-4-carboxylate
6630-33-7, 2-Bromobenzaldehyde 17259-77-7 17284-97-8 18908-07-1,
3-Methoxyphenylisocyanate 75680-93-2 78190-11-1,
1-[(Benzyloxy) carbonyl]piperidine-3-carboxylic acid 821002-63-5

821002-67-9 821002-71-5 RL: RCT (Reactant); RACT (Reactant or reagent)

(preparation of bone mass-increasing agents containing nonpeptidic osteoblast

differentiation promoting compds. and bisphosphonates) 448190-45-2P 596825-36-4P 596825-37-5P 596825-38-6P 596825-39-7P 596825-40-0P 596825-41-1P 596825-42-2P 596825-43-3P 596825-44-4P 596825-45-5P 596825-46-6P 596825-47-7P 596825-48-8P 596825-49-9P 596825-50-2P 596825-51-3P 596825-52-4P 596825-53-5P 596825-54-6P 596825-55-7P 596825-56-8P 596825-57-9P 596825-58-0P 596825-59-1P 596825-60-4P 596825-61-5P 596825-62-6P 596825-63-7P 596825-64-8P 596825-69-3P 596825-65-9P 596825-66-0P 596825-67-1P 596825-68-2P 596825-70-6P 596825-71-7P 596825-72-8P 596825-73-9P 596825-74-0P 596825-75-1P 596825-76-2P 596825-78-4P 596825-79-5P 596825-77-3P 596825-80-8P 596825-81-9P 596825-82-0P 596825-83-1P 596825-84-2P 596825-85-3P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of bone mass-increasing agents containing nonpeptidic osteoblast

differentiation promoting compds. and bisphosphonates)

MSTR 1

G1 = 13

ΗŊ----G2

```
G2 = alkyl <containing 1-6 C> (opt. substd. by (1-3) G3)
```

G3 = OH / CO2H

G4 = bond

Patent location: claim 2

Note: or pharmaceutically acceptable salts

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

=> d scan 12

L2 48 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN

IN Ethanamine, N,N-diethyl-2-[2-(2-phenyl-1,3-benzodioxol-2-yl)ethoxy]-, 2-hydroxy-1,2,3-propanetricarboxylate (1:1)

MF C21 H27 N O3 . C6 H8 O7

CM 1

CM 2

$$\begin{array}{c} {\rm CO_2H} \\ {\rm HO_2C-CH_2-C-CH_2-CO_2H} \\ {\rm OH} \end{array}$$

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

- L2 48 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
- ${\tt IN} \qquad 1-{\tt Naphthaleneheptanoic\ acid,\ 8-[(2,2-{\tt dimethyl-1-oxo-4-penten-1-yl)oxy]-1-penten-1-yl)oxy} = 1-{\tt Naphthaleneheptanoic\ acid,\ 8-[(2,2-{\tt dimethyl-1-oxo-4-penten-1-yl)oxy}]-1-penten-1-yl)oxy} = 1-{\tt Naphthaleneheptanoic\ acid,\ 8-[(2,2-{\tt dimethyl-1-oxo-4-penten-1-yl)oxy}} = 1-{\tt Naphthaleneheptanoic\ acid,\$
- 1, 2, 6, 7, 8, 8a-hexahydro- β , δ , 6-trihydroxy-2-methyl-
- MF C25 H38 O7

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

- L2 48 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
- IN 1-Naphthaleneheptanoic acid, 8-[(2S)-2-(2,3-dimethylphenoxy)-1-oxobutoxy]-1,2,6,7,8,8a-hexahydro- β ,8,6-trihydroxy-2-methyl-,(RR,8R,15,25,65,85,8aR)-

MF C30 H42 O8 CI COM

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

- L2 48 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
- IN 2-Propenoic acid, 2-methyl-, 2-(dimethylamino)ethyl ester,
- 2-hydroxy-1,2,3-propanetricarboxylate (1:1), polymer with 2-propenamide and 2-propenoic acid (9CI)
- MF (C8 H15 N 02 . C6 H8 07 . C3 H5 N O . C3 H4 O2)x
- CM 1

CM

CM 3

CM 4

$$\begin{array}{c|c} \circ & \operatorname{CH}_2 \\ \parallel & \parallel \\ \operatorname{Me}_2\operatorname{N} - \operatorname{CH}_2 - \operatorname{CH}_2 - \circ - \operatorname{C} - \operatorname{C} - \operatorname{Me} \end{array}$$

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

=> s l1 sss full

S L1 SSS FUL FILE=REGISTRY

FULL SEARCH INITIATED 11:14:50 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 906193 TO ITERATE

89.0% PROCESSED 806359 ITERATIONS

99.0% PROCESSED 897404 ITERATIONS

100.0% PROCESSED 906193 ITERATIONS SEARCH TIME: 00.00.38

L4 23608 SEA SSS FUL L1 1 FILES SEARCHED...

S L4 SSS FUL FILE=MARPAT

FULL SEARCH INITIATED 11:15:28 FILE 'MARPAT'

FULL SCREEN SEARCH COMPLETED - 223897 TO ITERATE

15.9%	PROCESSED	35678	ITERATIONS				3267	ANSWERS
25.8%	PROCESSED	57665	ITERATIONS				5016	ANSWERS
46.0%	PROCESSED	102956	ITERATIONS	(7	INCOMPLETE)	8240	ANSWERS
59.7%	PROCESSED	133776	ITERATIONS	(28	INCOMPLETE)	10943	ANSWERS
65.3%	PROCESSED	146162	ITERATIONS	(51	INCOMPLETE)	12247	ANSWERS
74.1%	PROCESSED	165836	ITERATIONS	(72	INCOMPLETE)	14000	ANSWERS
80.7%	PROCESSED	180774	ITERATIONS	(81	INCOMPLETE)	14853	ANSWERS
84.3%	PROCESSED	188711	ITERATIONS	(98	INCOMPLETE)	15642	ANSWERS
88.8%	PROCESSED	198742	ITERATIONS	(120	INCOMPLETE)	16645	ANSWERS
89.7%	PROCESSED	200838	ITERATIONS	(125	INCOMPLETE)	16845	ANSWERS
90.4%	PROCESSED	202314	ITERATIONS	(128	INCOMPLETE)	17008	ANSWERS
92.5%	PROCESSED	207170	ITERATIONS	(140	INCOMPLETE)	17512	ANSWERS
92.8%	PROCESSED	207863	ITERATIONS	(142	INCOMPLETE)	17583	ANSWERS
93.5%	PROCESSED	209418	ITERATIONS	(149	INCOMPLETE)	17756	ANSWERS
95.2%	PROCESSED	213165	ITERATIONS	(155	INCOMPLETE)	18161	ANSWERS

95.2% PROCESSED 213245 ITERATIONS (157 INCOMPLETE) 18175 ANSWERS

22673 ANSWERS

23591 ANSWERS

23608 ANSWERS

95.8%	PROCESSED	214435	ITERATIONS	(161	INCOMPLETE)	18294	ANSWERS
96.0%	PROCESSED	215008	ITERATIONS	(164	INCOMPLETE)	18346	ANSWERS
96.0%	PROCESSED	215008	ITERATIONS	(164	INCOMPLETE)	18346	ANSWERS
96.5%	PROCESSED	216146	ITERATIONS	(168	INCOMPLETE)	18473	ANSWERS
96.6%	PROCESSED	216379	ITERATIONS	(169	INCOMPLETE)	18493	ANSWERS
96.6%	PROCESSED	216379	ITERATIONS	(169	INCOMPLETE)	18493	ANSWERS
97.2%	PROCESSED	217555	ITERATIONS	(170	INCOMPLETE)	18599	ANSWERS
97.4%	PROCESSED	218078	ITERATIONS	(171	INCOMPLETE)	18665	ANSWERS
98.3%	PROCESSED	220098	ITERATIONS	(172	INCOMPLETE)	18797	ANSWERS
98.5%	PROCESSED	220647	ITERATIONS	(173	INCOMPLETE)	18849	ANSWERS
98.6%	PROCESSED	220655	ITERATIONS	(174	INCOMPLETE)	18850	ANSWERS
98.8%	PROCESSED	221208	ITERATIONS	(175	INCOMPLETE)	18912	ANSWERS
98.8%	PROCESSED	221208	ITERATIONS	(175	INCOMPLETE)	18912	ANSWERS
98.8%	PROCESSED	221221	ITERATIONS	(176	INCOMPLETE)	18914	ANSWERS
98.8%	PROCESSED	221221	ITERATIONS	(176	INCOMPLETE)	18914	ANSWERS
98.8%	PROCESSED	221221	ITERATIONS	(176	INCOMPLETE)	18914	ANSWERS
98.8%	PROCESSED	221221	ITERATIONS	(176	INCOMPLETE)	18914	ANSWERS
99.1%	PROCESSED	221827	ITERATIONS	(178	INCOMPLETE)	18971	ANSWERS
99.1%	PROCESSED	221827	ITERATIONS	(178	INCOMPLETE)	18971	ANSWERS
99.1%	PROCESSED	221827	ITERATIONS	(178	INCOMPLETE)	18971	ANSWERS
99.1%	PROCESSED	221827	ITERATIONS	(178	INCOMPLETE)	18971	ANSWERS
99.1%	PROCESSED	221827	ITERATIONS	(178	INCOMPLETE)	18971	ANSWERS
99.1%	PROCESSED	221827	ITERATIONS	(178	INCOMPLETE)	18971	ANSWERS
	PROCESSED -User Break		ITERATIONS	(178	INCOMPLETE)	18971	ANSWERS
			ITERATIONS	,	170	THEOMETERS	10070	ANSWERS
	PROCESSED ENDED BY USE		TIERATIONS	(1/9	INCOMPLETE)	183 15	ANDWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BROJECTED ITERATIONS: 223897 TO 223897
PROJECTED ANSWERS: 18972 TO 19551

L5 18972 SEA SSS FUL L1

SEARCH TIME: 00.11.32

¹ FILES SEARCHED...

(FILE 'CAPLUS, MARPAT, REGISTRY' ENTERED AT 11:04:57 ON 24 MAR 2009) DELETE HISTORY YES

DELETE SELECT YES STRUCTURE UPLOADED L2 48 S L1 SSS SAM FILE=REGISTRY 50 S L2 SSS SAM FILE-MARPAT

L4 23608 S L1 SSS FUL FILE=REGISTRY L5 18972 S L4 SSS FUL FILE=MARPAT

=> file caplus

L3

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 523.76 523.98

FILE 'CAPLUS' ENTERED AT 11:27:25 ON 24 MAR 2009 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

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FILE COVERS 1907 - 24 Mar 2009 VOL 150 ISS 13 FILE LAST UPDATED: 23 Mar 2009 (20090323/ED)

Caplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2008.

CAS Information Use Policies apply and are available at:

http://www.cas.org/legal/infopolicy.html

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 14 L6 159563 L4

=>	е	odor/ct			
E#		FREQUENCY	AT		TERM
E1		0	11		ODONUS/CT
E2		1	11		ODONUS NIGER/CT
E3		0	2	>	ODOR/CT
E4		33213	29		ODOR AND ODOROUS SUBSTANCES/CT
E5		0	4		ODOR AND ODOROUS SUBSTANCES (L) MUSK-LIKE/CT
E6		0	4		ODOR AND ODOROUS SUBSTANCES (L) ODORIZATION/CT
E7		0	2		ODOR AND ODOROUS SUBSTANCES (L) OFF-/CT
E8		0	7		ODOR AND ODOROUS SUBSTANCES (L) OFF-ODOR/CT
E9		0	2		ODOR MOL. STRUCTURE-BIOL. ACTIVITY RELATIONSHIP/CT
E1)	0	2		ODOR MOL. STRUCTURE-PROPERTY RELATIONSHIP/CT
E1:	1	0	2		ODOR RECEPTOR/CT
E1:	2	0	2		ODOR STRUCTURE-ACTIVITY RELATIONSHIP/CT

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=> e e3+a11
E1
            0
                --> Odor/CT
        33213 USE Odor and Odorous substances/CT
E.2
******** END *******
=> set expand continuous perm
SET COMMAND COMPLETED
=> e e2+a11
E3
         4097 BT1 Physical and chemical properties/CT
E4
        33213
                 --> Odor and Odorous substances/CT
                    HNTE Valid heading during volume 76 (1972) to present.
          450
                    OLD Odorous substances/CT
E6
         1757
                    OLD Odors/CT
E7
                    UF Aroma/CT
EΑ
                    UF Fragrance/CT
                    UF Fragrant substances/CT
E9
                    UF Odiferous compds./CT
E10
E11
                    UF Odor/CT
E12
                    UF
                       Osmophore/CT
E13
                    UF
                        Smell/CT
                    UF
E14
                        Smell (odor)/CT
                    RT Air fresheners/CT
E15
          978
E16
         6191
                    RT Cosmetics and personal care products/CT
E17
                   RT Deodorization/CT
         7775
                  RT Essences/CT
RT Flavor/CT
E18
          2595
E19
       25688
E20
        4692
                   RT Olfaction/CT
          626
E21
                   RT Olfactory system/CT
        13864
E22
                   RT Perfumes/CT
E23
       22247
                    RT
                        Volatile substances/CT
E24
                    RTCS &-Decalactone/CT
E25
                    RTCS y-Octalactone/CT
E26
                    RTCS 2-Methylisoborneol/CT
E27
                    RTCS Ethyl 2-methylbutanoate/CT
E28
                    RTCS Furaneol/CT
E29
                    RTCS Geosmin/CT
E30
                    RTCS Trans-2-Heptenal/CT
                    RTCS Trans-3-Hexen-1-o1/CT
****** END *******
=> s e4-e14
         33213 "ODOR AND ODOROUS SUBSTANCES"/CT
          450 "ODOROUS SUBSTANCES"/CT
          1757 ODORS/CT
            0 AROMA/CT
            0 FRAGRANCE/CT
            0 "FRAGRANT SUBSTANCES"/CT
            0 "ODIFEROUS COMPDS."/CT
            0 ODOR/CT
            0 OSMOPHORE/CT
            3 SMELL/CT
            0 "SMELL (ODOR)"/CT
L7
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              ODORS/CT OR AROMA/CT OR FRAGRANCE/CT OR "FRAGRANT SUBSTANCES"/CT
               OR "ODIFEROUS COMPDS."/CT OR ODOR/CT OR OSMOPHORE/CT OR SMELL/C
              T OR "SMELL (ODOR)"/CT)
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=> s 16 and 17

1239 L6 AND L7

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=> e sweat/ct
E# FREQUENCY AT TERM
E32
      0
                2 SWCV/CT
E33
          0
                1
                      SWE/CT
E34
        1566 14 --> SWEAT/CT
E35
               11 SWEAT (L) CONGENITAL SENSORY NEUROPATHY WITH ANHIDROSI
         0
                      S/CT
E36
                      SWEAT BANDS/CT
        1135 11
E37
                      SWEAT GLAND/CT
         0 11
E38
                      SWEAT GLAND (L) APOCRINE/CT
E39
          0 11
                      SWEAT GLAND (L) APOECCRINE/CT
E40
          0 10
0 12
                      SWEAT GLAND (L) DUCT/CT
E41
                      SWEAT GLAND (L) ECCRINE/CT
E42
          0 10
                      SWEAT GLAND (L) EPITHELIUM/CT
E43
         224
                2
                      SWEAT GLANDS/CT
=> e e34+a11
E44
              BT3 Biological processes and phenomena (non-CA heading)/CT
E45
           0
                BT2 Animal processes and phenomena (non-CA heading)/CT
                 BT2 Body, anatomical/CT
         8332
E46
E47
        23703
                  BT1 Body fluid/CT
                   BT1 Secretions (external)/CT
E48
         357
E49
         1566
                     --> Sweat/CT
                       HNTE Valid heading during volume 126 (1997) to
                           present.
E50
        2308
                       OLD Perspiration/CT
E51
                       UF Diaphoresis/CT
E52
                       UF Diaphoretics/CT
E53
                      UF Eccrine sweat/CT
        2414
E54
                      RT Antiperspirants/CT
E55
         291
                      RT Personal deodorants/CT
E56
      133490
                      RT Skin/CT
E57
        1135
                      RT Sweat gland/CT
****** END *******
=> s e49-e53
         1566 SWEAT/CT
         2308 PERSPIRATION/CT
            0 DIAPHORESIS/CT
            0 DIAPHORETICS/CT
            0 "ECCRINE SWEAT"/CT
L9
         3874 (SWEAT/CT OR PERSPIRATION/CT OR DIAPHORESIS/CT OR DIAPHORETICS/C
             T OR "ECCRINE SWEAT"/CT)
=> s 18 and 19
            8 I.8 AND I.9
T-10
=> d 1-8 ibib hit
L10 ANSWER 1 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 2007:269781 CAPLUS
DOCUMENT NUMBER:
                       146:280302
TITLE:
                       Biochemistry of human axilla malodor and chemistry of
                      deodorant ingredients
AUTHOR(S):
                      Gautschi, Markus; Natsch, Andreas; Schroeder, Fridtjof
CORPORATE SOURCE:
                      Fragrance Research, Givaudan Schweiz AG, Dubendorf,
                      CH-8600, Switz.
SOURCE:
                       Chimia (2007), 61(1-2), 27-32
                      CODEN: CHIMAD; ISSN: 0009-4293
PUBLISHER:
                      Swiss Chemical Society
DOCUMENT TYPE:
                      Journal; General Review
```

LANGUAGE: English

REFERENCE COUNT: 57 THERE ARE 57 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

Odor and Odorous substances

(axilla body malodor; human axilla malodor biochem, and chemical of

deodorant ingredients)

Sweat.

(axillary; human axilla malodor biochem. and chemical of deodorant ingredients)

1289-40-3, Hexenoic acid 58888-76-9, 3-Hydroxy-3-methyl-hexanoic

RL: BSU (Biological study, unclassified); BIOL (Biological study) (human axilla malodor biochem. and chemical of deodorant ingredients)

L10 ANSWER 2 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2006:1279662 CAPLUS

DOCUMENT NUMBER: 146.32449

TITLE: Oil-containing deodorizing aerosol compositions having

skin-cooling active substances

INVENTOR(S): Emmerling, Winfried; Heinsohn, Ulrike

PATENT ASSIGNEE(S): Henkel Kommanditgesellschaft Auf Aktien, Germany PCT Int. Appl., 20pp.

SOURCE: CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: German

FAMILY ACC. NUM. COUNT: 1 PATENT INFORMATION:

		TENT						DATE									ATE	
	WO	2006	1286	22		A2		2006	1207									
								AU,		BA.	BB.	BG.	BR.	BW.	BY.	B7.	CA.	CH.
								DE.										
								ID,										
								LT,										
								NZ,										
								TJ,										
								10,	111,	TIM,	IR,	11,	14,	UM,	uu,	05,	04,	vc,
		DIT.				ZM,		OF	DE.	DIZ		m.c	-	ED.	CD.	OD.	****	T 17
		RW:						CZ,										
								MC,										
								GN,										
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						RU,												
		1020																
	ΑU	2006	2543	59		A1		2006	1207		AU 2	006-:	2543.	59		2	0060.	524
	EΡ	1888	012			A2		2008	0220		EP 2	006-	7538:	30		2	0060.	524
		R:	AT,	BE,	BG,	CH,	CY,	CZ,	DE,	DK,	EE,	ES,	FI,	FR,	GB,	GR,	HU,	IE,
			IS,	IT,	LI,	LT,	LU,	LV,	MC,	NL,	PL,	PT,	RO,	SE,	SI,	SK,	TR	
	US	2008	0124	282		A1		2008	0529		US 2	007-	9481	92		2	0071	130
PRIO	RIT	Y APP	LN.	INFO	. :						DE 2	005-	1020	0502	5495	A 2	0050	601
											WO 2	006-1	EP49	31	1	W 2	0060	524
REFE	RENO	E CO	UNT:			3	Т	HERE	ARE									
		00				-		ECOD										

R HIS RECORD, ALL CITATIONS AVAILABLE IN THE RE FORMAT

Odor and Odorous substances (Protectate HR & MOD2;; oil-containing deodorizing aerosol compns. having skin-cooling active substances)

Sweat

P

(decomposing enzymes, inhibitors of; oil-containing deodorizing aerosol compns. having skin-cooling active substances)

50-21-5, Lactic acid, biological studies 65-85-0D, Benzoic acid, C12-C15 esters 69-72-7, Salicylic acid, biological studies 74-98-6, n-Propane,

```
biological studies 75-28-5, Isobutane 77-92-9, Citric acid,
     biological studies 78-78-4, Isopentane 79-14-1, Glycolic acid, biological studies 87-69-4, Tartaric acid, biological studies 89-78-1,
     Menthol 89-79-2, Isopulegol 103-23-1, Di(2-ethylhexyl)adipate
     104-76-7, 2-Ethylhexyl alcohol 105-99-7, Di-n-butyladipate 106-97-8,
     n-Butane, biological studies 109-43-3, Di-n-butylsebacate 109-66-0,
     n-Pentane, biological studies 110-27-0, Isopropyl myristate 110-40-7,
     Diethylsebacate 112-10-7, Isopropyl stearate 112-11-8, Isopropyloleate
     123-95-5, n-Butvlstearate 142-16-5, Dioctvlmaleate 142-91-6, Isopropvl
     palmitate 624-03-3, Ethyleneglycoldipalmitate 629-82-3, Cetiol OE
     693-19-6D, Isononanoic acid, cetearyl derivs. 928-24-5,
     Ethyleneglycoldioleate 1327-41-9, Reach 103 1565-76-0,
     Menthylmethylether 2432-87-3, Dioctylsebacate 2456-28-2, Didecyl ether
     2915-53-9, Dicapryl maleate 2915-57-3, Di-2-ethylhexylsuccinate
     3687-45-4, Oleyloleate 3687-46-5 4602-84-0, Farnesol 5444-75-7,
     Ethylhexylbenzoate 6915-15-7, Malic acid 6938-94-9,
     Diisopropyladipate 7491-02-3, Diisopropylsebacate
                                                            15763-02-7.
     Dioctylmalate 17071-54-4, Hexyl octyl ether 17088-93-6 17162-29-7, Menthyllactate 17618-45-0, 2-Hexyldecylstearate 17673-56-2,
     Oleylerucate 20292-08-4, 2-Ethylhexyllaurate 22047-49-0,
     2-Ethylhexylstearate 22766-84-3, 2-Octyldodecylpalmitate
                                                                    25339-09-7,
     Isocetylstearate 27215-38-9, Glycerin monolaurate 27640-89-7,
     Erucylerucate 28880-24-2, Diisooctylsuccinate, biological studies
     29806-73-3, 2-Ethylhexyl palmitate 30500-51-7, Isononvlstearate
     34316-64-8, n-Hexyllaurate 34364-24-4, Isostearylbenzoate Octyldodecanol 40550-16-1, Isooctylstearate 42131-25-9,
     Isononylisononanoate 42131-27-1, İsotridecylisononanoate 60209-82-7, Isodecylneopentanoate 63187-91-7, Menthone glycerin acetal 68171-33-5,
                                                                    60209-82-7.
     Isopropylisostearate 70445-33-9, Sensiva SC 50 74565-11-0, Finsolv TN
     75363-56-3 81897-25-8, 2-Ethylhexylisostearate 85617-81-8,
     Erucyloleate 92353-16-7, Hexyldecanol 117356-20-4 143608-26-8
     156324-82-2 156679-39-9 158599-25-8, Diglycerin monocaprinate
     161544-28-1, Cosmacol EMI
                                 195060-85-6 221333-66-0 227450-65-9,
     Hexvldecvllaurate 351420-48-9 841309-69-1
     RL: COS (Cosmetic use); BIOL (Biological study); USES (Uses)
        (oil-containing deodorizing aerosol compns. having skin-cooling active
        substances)
L10 ANSWER 3 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER:
                         2006:104126 CAPLUS
DOCUMENT NUMBER:
                          145:184938
TITLE:
                          A broad diversity of volatile carboxylic acids,
                          released by a bacterial aminoacylase from axilla
                          secretions, as candidate molecules for the
                          determination of human-body odor type
                          Natsch, Andreas; Derrer, Samuel; Flachsmann, Felix;
AUTHOR(S):
                          Schmid, Joachim
CORPORATE SOURCE:
                         Givaudan Schweiz AG, Duebendorf, CH-8600, Switz.
SOURCE:
                          Chemistry & Biodiversity (2006), 3(1), 1-20
                          CODEN: CBHIAM; ISSN: 1612-1872
                          Verlag Helvetica Chimica Acta AG
PUBLISHER:
DOCUMENT TYPE:
                          Journal
LANGUAGE:
                          English
REFERENCE COUNT:
                                THERE ARE 28 CITED REFERENCES AVAILABLE FOR THIS
```

Sweat (volatile carboxylic acids from axilla secretions, as candidate mols. for determination of human-body odor type)

(body odor; volatile carboxylic acids from axilla secretions, as candidate mols. for determination of human-body odor type)

Odor and Odorous substances

Human

RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

```
IT 50-21-5, Lactic acid, biological studies 103-82-2, Phenylacetic acid,
    biological studies 123-99-9, Azelaic acid, biological studies
    156-38-7, (4-Hydroxyphenyl)acetic acid 505-48-6, Suberic acid
    816-66-0 1460-34-0 3788-56-5, 9-Hydroxynonanoic acid 6966-34-3
    14292-26-3 14292-27-4 16493-80-4 27960-21-0 54068-86-9
    58888-76-9 59866-91-0 80113-38-8 80113-39-9
    132735-95-6 160595-71-1 875712-96-2
                                           903503-32-2
    903503-33-3 903503-34-4 903503-35-5
    RL: BSU (Biological study, unclassified); BIOL (Biological study)
        (volatile carboxylic acids from axilla secretions, as candidate mols.
       for determination of human-body odor type)
L10 ANSWER 4 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER:
                        2005:329511 CAPLUS
DOCUMENT NUMBER:
                        143:332012
TITLE:
                        Isolation of a bacterial enzyme releasing axillary
                        malodor and its use as a screening target for novel
                        deodorant formulations
AUTHOR(S):
                        Natsch, A.; Gfeller, H.; Gygax, P.; Schmid, J.
CORPORATE SOURCE:
                        Givaudan Schweiz AG, Duebendorf, CH-8600, Switz.
SOURCE:
                        International Journal of Cosmetic Science (2005),
                        27(2), 115-122
                        CODEN: IJCMDW; ISSN: 0142-5463
PUBLISHER:
                        Blackwell Publishing Ltd.
DOCUMENT TYPE:
                        Journal
LANGUAGE:
                        English
REFERENCE COUNT:
                              THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS
                              RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT
    Sweat
        (axillary; isolation of bacterial enzyme releasing axillary malodor)
    Odor and Odorous substances
        (off-odor; isolation of bacterial enzyme releasing axillary malodor)
    58888-76-9P, 3-Hydroxy-3-methyl-hexanoic acid
    RL: ANT (Analyte); PUR (Purification or recovery); ANST (Analytical
    study); PREP (Preparation)
        (isolation of bacterial enzyme releasing axillary malodor)
L10 ANSWER 5 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER:
                        2005:19524 CAPLUS
DOCUMENT NUMBER:
                        142:313793
                        Identification of new odoriferous compounds in human
TITLE:
                        axillary sweat
AUTHOR(S):
                        Hasegawa, Yoshihiro; Yabuki, Masayuki; Matsukane,
                        Masamoto
CORPORATE SOURCE:
                        Perfumery Development Research Laboratory, Kao
                        Corporation, Tokyo, Japan
                        Chemistry & Biodiversity (2004), 1(12), 2042-2050
SOURCE:
                        CODEN: CBHIAM; ISSN: 1612-1872
PUBLISHER:
                        Verlag Helvetica Chimica Acta AG
DOCUMENT TYPE:
                        Journal
LANGUAGE:
                        English
REFERENCE COUNT:
                        19
                              THERE ARE 19 CITED REFERENCES AVAILABLE FOR THIS
                              RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT
    Human
      Odor and Odorous substances
        (identification of new odoriferous compds. in human axillary sweat)
    51755-83-0P 58888-76-9P, 3-Hydroxy-3-methylhexanoic acid
    227456-33-9P 307964-33-6P
                                 548740-99-4P 757219-24-2P
    757219-38-8P 828300-39-6P
    RL: BSU (Biological study, unclassified); PRP (Properties); SPN (Synthetic
```

preparation); BIOL (Biological study); PREP (Preparation)

L10 ANSWER 6 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2004:754422 CAPLUS

DOCUMENT NUMBER: 141:282457

TITLE: Pseudo-body odor composition and perfume composition

for inhibiting body odor

INVENTOR(S): Ogura, Miharu; Sakurai, Kazutoshi; Sawano, Kiyohito;

Yamazaki, Sadahiko; Hirano, Koji

PATENT ASSIGNEE(S): Takasago International Corporation, Japan

SOURCE: PCT Int. Appl., 111 pp.

CODEN: PIXXD2
DOCUMENT TYPE: Patent

LANGUAGE: Facent English

FAMILY ACC. NUM. COUNT: 1

PA'	TENT	NO.			KIN	D	DATE			APPI	LICAT	ION 1	NO.		D.	ATE	
WO	2004	0781	54		A1		2004	0916		WO 2	2004-	JP23	00		2	0040	226
	W:	ΑE,	AG,	AL,	AM,	ΑT,	AU,	AZ,	BA,	BB,	BG,	BR,	BW,	BY,	ΒZ,	CA,	CH,
		CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD,
		GE,	GH,	GM,	HR,	ΗU,	ID,	IL,	IN,	IS,	KE,	KG,	KP,	KR,	ΚZ,	LC,	LK,
		LR,	LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	ΜZ,	NA,	NΙ,	NO
	RW:	BW,	GH,	GM,	KE,	LS,	MW,	ΜZ,	SD,	SL,	SZ,	TZ,	UG,	ZM,	ZW,	ΑT,	BE,
		BG,	CH,	CY,	CZ,	DE,	DK,	EE,	ES,	FI,	FR,	GB,	GR,	HU,	ΙE,	IT,	LU,
		MC,	NL,	PT,	RO,	SE,	SI,	SK,	TR,	BF,	BJ,	CF,	CG,	CI,	CM,	GΑ,	GN,
		GQ,	GW,	ML,	MR,			TD,									
	2004										2003-						
	2004																
AU	2004	2185	60		A1		2004	0916		AU 2	2004-	2185	60		2	0040	226
EP	1601																
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											TR,						
	1756									CN 2	2004-	8000	5817		2	0040	226
	1003						2007										
	2006				A1		2006	0720			2005-						
PRIORIT	Y APP	LN.	INFO	. :							2003-					0030	
											2003-					0030	
										WO 2	2004-	JP23	00	- 1	A 2	0040	226

REFERENCE COUNT: 8 THERE

THERE ARE 8 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

IT Odor and Odorous substances

(body, pseudo-, for deodorant evaluation; pseudo-body odor composition containing aldehydes, carboxylic acids and fatty acids for evaluation of deodorant perfumes for inhibiting body odor)

IT Sweat

(pseudo-, for deodorant evaluation; pseudo-body odor composition containing aldehydes, carboxylic acids and fatty acids for evaluation of deodorant perfumes for inhibiting body odor)

IT 50-21-5, Lactic acid, biological studies 64-19-7, Acetic acid, biological studies 77-93-0, Triethyl citrate 79-31-2, Isobutanoic acid 107-92-6, Butanoic acid, biological studies 111-14-8, Heptanoic acid 112-05-0, Nonanoic acid 112-44-7, Undecyl aldehyde 120-51-4, Benzyl benzoate 124-07-2, Octanoic acid, biological studies 142-62-1, Hexanoic acid, biological studies 503-74-2, Isovaleric acid 27960-21-0, (E)-3-Methyl-2-hexanoic acid 54068-86-9, (2)-3-Methyl-2-hexanoic acid 58088-76-9, 3-Hydroxy-3-methylhexanoic acid 557219-24-2 757219-38-8 RL: BUU (Biological use, unclassified); BIOL (Biological study); USES

(Uses)
(pseudo-body odor composition containing aldehydes, carboxylic acids and

L10 ANSWER 7 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2004:308604 CAPLUS

DOCUMENT NUMBER: 140:326641

TITLE: Indicator for assessing body odor, process for producing the same, body odor assessment method,

method of assessing efficaciousness of deodorant and

kit for conveniently assessing body odor INVENTOR(S):

Yabuki, Masavuki; Hasegawa, Yoshihiro; Matsukane,

Masamoto; Yabe, Emi

PATENT ASSIGNEE(S): Kao Corporation, Japan SOURCE: PCT Int. Appl., 86 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

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V												2003-					0031	006
		W:	ΑE,	AG,	AL,	AM,	AT,	AU,	AZ,	BA,	BB	, BG,	BR,	BY,	BZ,	CA,	CH,	CN,
			CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC	, EE,	EG,	ES,	FΙ,	GB,	GD,	GE,
			GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	KE	, KG,	KP,	KR,	KZ,	LC,	LK,	LR,
												, MW,						
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		R:										, IT,						
												TR,						
											JP :	2004-	8365	4		2	0040	322
	JΡ	4081	034			B2		2008	0423							_		
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												2003-				A 2		
												2003-						
											WO :	2003-	JP12	193	1	N 2	0031	UU6

OTHER SOURCE(S): MARPAT 140:326641 REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS

RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

Deodorants (personal)

Human

Odor and Odorous substances

Sweat

(indicator for assessing body odor containing \(\beta\)-hydroxy carboxylates or 3-mercapto alcs., process for producing same, body odor assessment method, method of assessing efficaciousness of deodorant and kit for conveniently assessing body odor)

51755-83-0P, 3-Mercaptohexanol 58888-76-9P,

3-Hydroxy-3-methylhexanoic acid 227456-27-1P 227456-33-9P

307964-23-4P 548740-99-4P

RL: ANT (Analyte); BSU (Biological study, unclassified); PUR (Purification

or recovery); ANST (Analytical study); BIOL (Biological study); PREP (Preparation)

(indicator for assessing body odor containing β-hydroxy carboxylates or 3-mercapto alcs., process for producing same, body odor assessment method, method of assessing efficaciousness of deodorant and kit for conveniently assessing body odor)

L10 ANSWER 8 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2003:122930 CAPLUS

DOCUMENT NUMBER: 139:2719

TITLE: A Specific Bacterial Aminoacylase Cleaves Odorant

Precursors Secreted in the Human Axilla

AUTHOR(S): Natsch, Andreas; Gfeller, Hans; Gygax, Peter; Schmid, Joachim; Acuna, Gonzalo

Givaudan Duebendorf Ltd., Duebendorf, CH-8600, Switz. CORPORATE SOURCE:

SOURCE: Journal of Biological Chemistry (2003), 278(8),

5718-5727

CODEN: JBCHA3; ISSN: 0021-9258

PUBLISHER: American Society for Biochemistry and Molecular

Biology DOCUMENT TYPE: Journal

LANGUAGE: English

REFERENCE COUNT: THERE ARE 33 CITED REFERENCES AVAILABLE FOR THIS 33 RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

Odor and Odorous substances

(anal. of hydrolyzed axilla secretions)

35205-70-0, 3-Methyl-2-hexenoic acid 58888-76-9,

3-Hvdroxv-3-methvlhexanoic acid

RL: BSU (Biological study, unclassified); BIOL (Biological study) (anal. of hydrolyzed axilla secretions)

=> FIL REGISTRY

COST IN U.S. DOLLARS SINCE FILE

TOTAL ENTRY SESSION FULL ESTIMATED COST 52.88 576.86

FILE 'REGISTRY' ENTERED AT 11:32:28 ON 24 MAR 2009

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

22 MAR 2009 HIGHEST RN 1125392-64-4 STRUCTURE FILE UPDATES: DICTIONARY FILE UPDATES: 22 MAR 2009 HIGHEST RN 1125392-64-4

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 9, 2009.

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/support/stngen/stndoc/properties.html

=> S 58888-76-9/RN

L11 1 58888-76-9/RN

=> SET NOTICE 1 DISPLAY

NOTICE SET TO 1 U.S. DOLLAR FOR DISPLAY COMMAND SET COMMAND COMPLETED

=> D L11 SOIDE 1-

YOU HAVE REQUESTED DATA FROM 1 ANSWERS - CONTINUE? Y/(N):v THE ESTIMATED COST FOR THIS REQUEST IS 6.85 U.S. DOLLARS DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y) /N:y

- L11 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2009 ACS on STN RN 58888-76-9 REGISTRY
- CN Hexanoic acid, 3-hydroxy-3-methyl- (CA INDEX NAME)
- OTHER NAMES:

CN

- β-Hydroxy-β-methylcaproic acid CN β-Hydroxy-β-methylhexanoic acid
- CN 3-Hydroxy-3-methylhexanoic acid
- MF C7 H14 O3
- STN Files: BEILSTEIN*, BIOSIS, CA, CAPLUS, CHEMCATS, TOXCENTER, USPATFULL.

(*File contains numerically searchable property data)

- DT.CA CAplus document type: Journal; Patent RL.P
- Roles from patents: ANST (Analytical study); BIOL (Biological study); PREP (Preparation); USES (Uses)
- RL.NP Roles from non-patents: ANST (Analytical study); BIOL (Biological study); PREP (Preparation); PRP (Properties); USES (Uses)

- **PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**
 - 15 REFERENCES IN FILE CA (1907 TO DATE) 15 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> SET NOTICE LOGIN DISPLAY

NOTICE SET TO OFF FOR DISPLAY COMMAND SET COMMAND COMPLETED

=> d ibib abs hitstr 1-8 'IBIB' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

'ABS' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG RN

SAM - Index Name, MF, and structure - no RN

FIDE - All substance data, except sequence data

- FIDE, but only 50 names

SOIDE - IDE, plus sequence data

SOIDE3 - Same as SOIDE, but 3-letter amino acid codes are used

SOD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used

SON - Protein sequence name information, includes RN

EPROP - Table of experimental properties

PPROP - Table of predicted properties

PROP - EPROP, ETAG, PPROP and SPEC

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract

APPS -- Application and Priority Information

BIB -- CA Accession Number, plus Bibliographic Data

CAN -- CA Accession Number

CBIB -- CA Accession Number, plus Bibliographic Data (compressed)

IND -- Index Data

IPC -- International Patent Classification

PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS -- ABS, indented, with text labels

IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)

OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations

SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields. HELP FORMATS -- To see detailed descriptions of the predefined formats. ENTER DISPLAY FORMAT (IDE):wend

'WEND' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

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REG - RN
SAM
      - Index Name, MF, and structure - no RN
FIDE - All substance data, except sequence data
IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data
SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used
SQD
    - Protein sequence data, includes RN
SOD3 - Same as SOD, but 3-letter amino acid codes are used
SON
    - Protein sequence name information, includes RN
EPROP - Table of experimental properties
PPROP - Table of predicted properties
PROP - EPROP, ETAG, PPROP and SPEC
Any CA File format may be combined with any substance format to
obtain CA references citing the substance. The substance formats
must be cited first. The CA File predefined formats are:
ABS -- Abstract
APPS -- Application and Priority Information
BIB -- CA Accession Number, plus Bibliographic Data
CAN -- CA Accession Number
CBIB -- CA Accession Number, plus Bibliographic Data (compressed)
IND -- Index Data
IPC -- International Patent Classification
PATS -- PI, SO
STD -- BIB, IPC, and NCL
IABS -- ABS, indented, with text labels
IBIB -- BIB, indented, with text labels
ISTD -- STD format, indented
OBIB ----- AN, plus Bibliographic Data (original)
OIBIB ----- OBIB, indented with text labels
SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations
The ALL format gives FIDE BIB ABS IND RE, plus sequence data when
it is available.
The MAX format is the same as ALL.
The IALL format is the same as ALL with BIB ABS and IND indented.
with text labels.
For additional information, please consult the following help
messages:
HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):end
=> file caplus
COST IN U.S. DOLLARS
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                                                              TOTAL
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FILE 'CAPLUS' ENTERED AT 11:33:18 ON 24 MAR 2009
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PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
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FULL ESTIMATED COST

ENTRY

2.53

SESSION

579.39

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FILE COVERS 1907 - 24 Mar 2009 VOL 150 ISS 13
FILE LAST UPDATED: 23 Mar 2009 (20090323/ED)
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Caplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2008.

CAS Information Use Policies apply and are available at:

http://www.cas.org/legal/infopolicy.html

This file contains CAS Registry Numbers for easy and accurate substance identification.

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=> display ibib abs hitstr 1-8
ENTER (L11), L# OR ?:y
'Y' IS NOT VALID HERE
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For an explanation, enter "HELP DISPLAY".

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=> 111
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L11 IS NOT A RECOGNIZED COMMAND

3874 S E49-E53

8 S L8 AND L9

The previous command name entered was not recognized by the system. For a list of commands available to you in the current file, enter "HELP COMMANDS" at an arrow prompt (=>).

=> display ibib abs hitstr 1-8 111
YOU HAVE REQUESTED DATA FROM FILE 'REGISTRY' - CONTINUE? (Y)/N:end

DELETE HISTORY YES

=> d his

T. 9

L10

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DELETE SELECT YES
                STRUCTURE UPLOADED
1.2
             48 S L1 SSS SAM FILE=REGISTRY
1.3
             50 S L2 SSS SAM FILE=MARPAT
          23608 S L1 SSS FUL FILE=REGISTRY
L4
L5
          18972 S L4 SSS FUL FILE=MARPAT
     FILE 'CAPLUS' ENTERED AT 11:27:25 ON 24 MAR 2009
L6
         159563 S L4
                E ODOR/CT
                E E3+ALL
                SET EXPAND CONTINUOUS PERM
                E E2+ALL
          35413 S E4-E14
1.8
           1239 S L6 AND L7
                E SWEAT/CT
                E E34+ALL
```

(FILE 'CAPLUS, MARPAT, REGISTRY' ENTERED AT 11:04:57 ON 24 MAR 2009)

FILE 'REGISTRY' ENTERED AT 11:32:28 ON 24 MAR 2009 1 S 58888-76-9/RN

SET NOTICE 1 DISPLAY SET NOTICE LOGIN DISPLAY

FILE 'CAPLUS' ENTERED AT 11:33:18 ON 24 MAR 2009

=> d 110 1-8 ibib abs hitstr

L10 ANSWER 1 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2007:269781 CAPLUS

DOCUMENT NUMBER: 146:280302

TITLE: Biochemistry of human axilla malodor and chemistry of

deodorant ingredients

AUTHOR(S): Gautschi, Markus; Natsch, Andreas; Schroeder, Fridtjof CORPORATE SOURCE: Fragrance Research, Givaudan Schweiz AG, Dubendorf,

CH-8600, Switz.

SOURCE: Chimia (2007), 61(1-2), 27-32 CODEN: CHIMAD; ISSN: 0009-4293

PUBLISHER: Swiss Chemical Society
DOCUMENT TYPE: Journal; General Review

LANGUAGE: English

A review. Key human axilla malodorants are hexenoic acid and 3-hydroxy-3-methyl-hexanoic acid, accompanied by some 25 structurally related hydroxyalkanoic acids and alkenoic acids. These sweat acids are secreted in the axilla in the form of odorless glutamine conjugates and are released upon enzymolysis by AMRE (axillary malodor releasing enzyme), produced by Corynebacteria. The sulfanylalkanols represent another important group of axilla malodor compds. that are also secreted in the form of odorless precursors. The major precursors are cysteine-glycine sulfanylalkanol conjugates and the minor precursors are cysteine sulfanylalkanol conjugates. The release occurs upon action of a β -lyase of axilla bacteria. Besides the classical approaches of axilla malodor masking using fragrances, the use of chems. to neutralize malodorants is described. The elucidation of the biochem. of the sweat acid release has allowed the development of fragrance precursors that act as competitive substrates to the natural malodor precursors as well as the development of specific antagonists that block AMRE. Finally, the characterization and functional expression of a first human malodor receptor presents an interesting approach for future development of axilla malodor blockers.

IT 58888-76-9, 3-Hydroxy-3-methyl-hexanoic acid

RL: BSU (Biological study, unclassified); BIOL (Biological study) (human axilla malodor biochem. and chemical of deodorant ingredients)

RN 58888-76-9 CAPLUS

CN Hexanoic acid, 3-hydroxy-3-methyl- (CA INDEX NAME)

REFERENCE COUNT: 57 THERE ARE 57 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L10 ANSWER 2 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN ACCESSION NUMBER: 2006:1279662 CAPLUS DOCUMENT NUMBER: 146:32449

TITLE:

Oil-containing deodorizing aerosol compositions having

skin-cooling active substances

INVENTOR(S): PATENT ASSIGNEE(S): Emmerling, Winfried; Heinsohn, Ulrike

Henkel Kommanditgesellschaft Auf Aktien, Germany

SOURCE: PCT Int. Appl., 20pp. CODEN: PIXXD2

DOCUMENT TYPE: Pat.ent. LANGUAGE: German

FAMILY ACC. NUM. COUNT: 1 PATENT INFORMATION:

		ENT I						DATE									ATE	
	WO	2006	1286	22		A2		2006	1207			006-					0060	
	WU																	
		W:	ΑE,	AG,	AL,	ΑM,	AT,	AU,	ΑZ,	BA,	BB,	BG,	BR,	BW,	BY,	ΒZ,	CA,	CH,
			CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD,
			GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	KM,	KN,	KP,	KR,
			KZ.	LC.	LK.	LR.	LS.	LT.	LU.	LV.	LY.	MA.	MD.	MG.	MK.	MN.	MW.	MX.
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											WO 2	006-	EP49	31	1	W 2	0060	524

- ΔR The invention relates to essentially alc.-free deodorizing aerosol compns., which contain selected oils, a skin-cooling active substance, a deodorizing active substance and a propellant gas. Thus a composition contained (weight/weight%): iso-Pr myristate 5.0; tri-Et citrate 2.9; phenoxyethanol 0.5; Proctate HR 0.2; α-(2-ethylhexyl)glycerin ether 0.3; diglycerin monocaprinate 0.2; Optacool A (not encapsulated) 0.1; perfume (not encapsulated) 1.0; n-butane to 100.
- 77-92-9, Citric acid, biological studies 6915-15-7, Malic acid

RL: COS (Cosmetic use); BIOL (Biological study); USES (Uses)

(oil-containing deodorizing aerosol compns. having skin-cooling active substances)

77-92-9 CAPLUS RN

1,2,3-Propanetricarboxylic acid, 2-hydroxy- (CA INDEX NAME) CN

RN 6915-15-7 CAPLUS

CN Butanedioic acid, 2-hydroxy- (CA INDEX NAME)

OH HO2C-CH-CH2-CO2H

REFERENCE COUNT:

3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD, ALL CITATIONS AVAILABLE IN THE RE FORMAT

L10 ANSWER 3 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2006:104126 CAPLUS

DOCUMENT NUMBER: 145:184938

TITLE: A broad diversity of volatile carboxylic acids, released by a bacterial aminoacylase from axilla secretions, as candidate molecules for the

determination of human-body odor type Natsch, Andreas; Derrer, Samuel; Flachsmann, Felix; AUTHOR(S):

Schmid, Joachim

CORPORATE SOURCE: Givaudan Schweiz AG, Duebendorf, CH-8600, Switz.

SOURCE: Chemistry & Biodiversity (2006), 3(1), 1-20 CODEN: CBHIAM; ISSN: 1612-1872 PUBLISHER: Verlag Helvetica Chimica Acta AG

DOCUMENT TYPE: Journal

LANGUAGE: English

Human body odor is to a large part determined by secretions of glands in the axillary regions. Two key odoriferous principles, 3-methylhex-2-enoic acid (3MH2; 4/5) and 3-hydroxy-3-methylhexanoic acid (HMHA; 6) have been shown to be released from glutamine conjugates secreted in the axilla by a specific Na-acyl-glutamine aminoacylase (N-AGA) obtained from axilla isolates of Corynebacteria sp. However, the low number of different odorants reported in humans stands in contrast to the observed high inter-individual variability in body odors. Axilla secretions of individual donors were, therefore, analyzed in detail. The secretions were treated with N-AGA, analyzed by GC/MS, and compared to undigested controls. Over 28 different carboxylic acids were released by this enzyme from odorless axilla secretions (Table 1). Many of these body odorants have not been reported before from a natural source, and they include several aliphatic 3-hydroxy acids with 4-Me branches, 3,4-unsatd., 4-Et-branched aliphatic acids, and a variety of degradation products of amino acids. The odor threshold of some of the acids was found to be in the range of 1 ng. Most of these compds. were present in all donors tested, but in highly variable relative amts., and they are, thus, candidate mols, as key components of a 'compound odor' determining the individual types of human body odor.

58888-76-9 59866-91-0 160595-71-1 RL: BSU (Biological study, unclassified); BIOL (Biological study) (volatile carboxylic acids from axilla secretions, as candidate mols. for determination of human-body odor type)

58888-76-9 CAPLUS RN

Hexanoic acid, 3-hvdroxv-3-methvl- (CA INDEX NAME) CN

RN 59866-91-0 CAPLUS

Hexanoic acid, 3-hydroxy-4-methyl- (CA INDEX NAME)

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

RN 160595-71-1 CAPLUS

REFERENCE COUNT:

28 THERE ARE 28 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L10 ANSWER 4 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN ACCESSION NUMBER: 2005:329511 CAPLUS

DOCUMENT NUMBER:

TITLE:

AUTHOR(S): CORPORATE SOURCE:

SOURCE:

PUBLISHER: DOCUMENT TYPE: 143:332012

Isolation of a bacterial enzyme releasing axillary malodor and its use as a screening target for novel deodorant formulations

Natsch, A.; Gfeller, H.; Gygax, P.; Schmid, J. Givaudan Schweiz AG, Duebendorf, CH-8600, Switz. International Journal of Cosmetic Science (2005), 27(2), 115-122

CODEN: IJCMDW; ISSN: 0142-5463 Blackwell Publishing Ltd.

Journal LANGUAGE: English

Axillary odor is known since 50 years to be formed upon the action of Corvnebacteria on odorless axilla secretions, but the nature of the bacterial enzymes involved in this process remained a mystery. The authors identified the known axilla odor determinant 3-methyl-2-hexenoic acid in hydrolyzed axilla secretions along with a new, chemical related compound, 3-hydroxy-3-methyl-hexanoic acid. The natural, odorless precursors of both these acids were purified from non-hydrolyzed fresh axilla secretions. The malodorous acids were shown to be covalently linked to a glutamine residue in fresh axilla secretions. Corynebacteria, but not Staphylococci, isolated from the axilla were found to release the acids from these precursors in vitro. A Zn2+-dependent aminoacylase mediating this cleavage was then purified from Corynebacterium striatum Ax20 and the corresponding gene agaA was cloned and heterologously expressed in Escherichia coli. Based on these biochem. findings, novel approaches in research on axilla malodor control are presented: (a) With a new test method using the isolated Corynebacteria and their enzymic activity, the direct malodor-controlling activity of existing cosmetic ingredients was evaluated. (b) The structure of the natural malodor precursor was modified by replacing the malodor acid with fragrance mols. These new fragrance precursors were shown to be cleaved by the same aminoacvlase.

58888-76-9P, 3-Hydroxy-3-methyl-hexanoic acid TТ RL: ANT (Analyte); PUR (Purification or recovery); ANST (Analytical study); PREP (Preparation) (isolation of bacterial enzyme releasing axillary malodor)

RN 58888-76-9 CAPLUS

Hexanoic acid, 3-hydroxy-3-methyl- (CA INDEX NAME) CN

REFERENCE COUNT: 7

THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L10 ANSWER 5 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2005:19524 CAPLUS

DOCUMENT NUMBER: 142:313793

Identification of new odoriferous compounds in human TITLE:

axillary sweat

AUTHOR(S): Hasegawa, Yoshihiro; Yabuki, Masavuki; Matsukane,

Masamoto

CORPORATE SOURCE: Perfumery Development Research Laboratory, Kao

Corporation, Tokyo, Japan

SOURCE: Chemistry & Biodiversity (2004), 1(12), 2042-2050

CODEN: CBHIAM; ISSN: 1612-1872

PUBLISHER: Verlag Helvetica Chimica Acta AG

DOCUMENT TYPE: Journal

LANGUAGE: English

3-Hydroxy-3-methylhexanoic acid (1) and the 3-sulfanylalkan-1-ols (2-5) were identified to contribute to the odor of human axillary sweat. Quant. analyses of axillary sweat exts. from 50 healthy men showed an unambiguous correlation between the detected levels of 1 and the intensity of the axillary odor. Chiral-GC analyses revealed 1 to be a 72: 28 mixture of the (S)/(R)-isomers. Optically pure (S)-1 (> 97% ee) emanated a strong spicy note, which recalled typical axillary odors. 3-Methyl-3-sulfanylhexan-1-ol (2), the enantiomeric ratio of which equaled that of 1, was present in greater quantity than any of the other

3-sulfanylalkanols. Optically pure (S)-2 (> 97% ee) had a strong meaty, fruity note, also reminiscent of axillary odor. The compds. identified, in particular (S)-1 and (S)-2, contribute significantly to the olfactory impression of human axillary odor.

58888-76-9P, 3-Hydroxy-3-methylhexanoic acid 757219-24-2P

757219-38-8P

RL: BSU (Biological study, unclassified); PRP (Properties); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)

(identification of new odoriferous compds. in human axillary sweat)

RN 58888-76-9 CAPLUS

CN Hexanoic acid, 3-hydroxy-3-methyl- (CA INDEX NAME)

$$\begin{array}{c} \text{OH} \\ \mid \\ \text{n-Pr-C-CH}_2\text{--CO}_2\text{H} \\ \mid \\ \text{Me} \end{array}$$

757219-24-2 CAPLUS RN

Hexanoic acid, 3-hvdroxv-3-methvl-, (3R)- (CA INDEX NAME) CN

Absolute stereochemistry. Rotation (-).

RN 757219-38-8 CAPLUS

Hexanoic acid, 3-hvdroxv-3-methyl-, (3S)- (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

REFERENCE COUNT: 19 THERE ARE 19 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L10 ANSWER 6 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2004:754422 CAPLUS

DOCUMENT NUMBER: 141:282457

TITLE: Pseudo-body odor composition and perfume composition for inhibiting body odor

INVENTOR(S): Ogura, Miharu; Sakurai, Kazutoshi; Sawano, Kiyohito;

Yamazaki, Sadahiko, Hirano, Koji

PATENT ASSIGNEE(S): Takasago International Corporation, Japan SOURCE: PCT Int. Appl., 111 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1 PATENT INFORMATION:

		ENT :										LICAT					ATE	
												2004-					0040	226
		W:	ΑE,	AG,	AL,	AM,	AT,	AU,	AZ,	BA,	BB,	, BG,	BR,	BW,	BY,	BZ,	CA,	CH,
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		0001							TD,		TD /	2003-		-		_	0020	202
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			IE,	SI,	LT,	LV,	FI,	RO,	MK,	CY,	AL,	, TR,	BG,	CZ,	EE,	HU,	SK	
	CN	1756	525			A		2006	0405		CN 2	2004-	8000	5817		2	0040	226
	CN	1003	5690	1		С		2007	1226									
	US	2006	0159	639		A1		2006	0720		US :	2005-	5476	38		2	0050	901
PRIOR	RIORITY APPLN. INFO.:										JP 2	2003-	5601	7	- 1	A 2	0030	303
											JP :	2003-	5746	2		A 2	0030	304
											wo :	2004-	JP23	00		A 2	0040	226

A pseudo-body odor composition according to the present invention comprises (A) AB at least one compound selected from hydroxyalkynyl carboxylic acids having 5 to 8 carbon atoms and alkenyl carboxylic acids having 5 to 8 carbon atoms and one double bond; (B) at least one compound selected from fatty acids having 2 to 24 carbon atoms; and (C) at least one compound selected from aldehydes having 2 to 13 carbon atoms. A deodorant perfume composition according to the invention includes one or more components that are excellent in masking effect and/or a harmonizing effect to an acid odor selected from hydroxyalkynyl carboxylic acids having 5 to 8 carbon atoms and alkenyl carboxylic acids having 5 to 8 carbon atoms and one double bond. For example, a composition that obviously senses the body odor likeliness, in particular as a sweat odor and armpit odor, contained 3-hydroxy-3-methylhexanoic acid (enantiomer ratio R/S of 1:3) 13.00, 3-methyl-2-hexenoic acid (trans/cis isomers ratio of 1:1) 5.00, acetic acid 11.85, isobutanoic acid 3.95, butanoic acid 0.79, isovaleric acid

- 3.95, hexanoic acid 0.79, heptanoic acid 0.79, octanoic acid 0.79, nonanoic acid 0.79, lactic acid 55.30, hexyl aldehyde 0.60, octyl aldehyde 0.36, nonyl aldehyde 0.99, evely aldehyde 0.90, undecyl aldehyde 0.60, dodecyl aldehyde 0.23, benzyl benzoate 0.23, and tri-Et citrate 5.00%, resp. Perfume compns. were also given and evaluated for their masking or harmonizing effects using a 2 cm x 2 cm filter paper with 200 μL of a pseudo body odor composition obtained.
- IT 58888-76-9, 3-Hydroxy-3-methylhexanoic acid 757219-24-2 757219-38-8

RL: BUU (Biological use, unclassified); BIOL (Biological study); USES

(pseudo-body odor composition containing aldehydes, carboxylic acids and fatty

acids for evaluation of deodorant perfumes for inhibiting body odor) RN 58888-76-9 CAPLUS

CN Hexanoic acid, 3-hydroxy-3-methyl- (CA INDEX NAME)

RN 757219-24-2 CAPLUS

N Hexanoic acid, 3-hydroxy-3-methyl-, (3R)- (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 757219-38-8 CAPLUS

CN Hexanoic acid, 3-hydroxy-3-methyl-, (3S)- (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

REFERENCE COUNT: 8 THERE ARE 8 CITED REFERENCES AVAILABLE FOR THIS
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L10 ANSWER 7 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2004:308604 CAPLUS DOCUMENT NUMBER: 140:326641

TITLE: Indicator for assessing body odor, process for producing the same, body odor assessment method,

method of assessing efficaciousness of deodorant and kit for conveniently assessing body odor

INVENTOR(S): Yabuki, Masayuki; Hasegawa, Yoshihiro; Matsukane,

PATENT ASSIGNEE(S): Kao Corporation, Japan SOURCE: PCT Int. Appl., 86 pp.

PCT Int. Appl., 86 pp. CODEN: PIXXD2 DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

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												2003-						
												BG,						
												, EE,						
												, KG,						
			LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN	, MW,	MX,	MZ,	NI,	NO.	NZ,	OM,
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			TR,	TT,	TZ,	UA,	UG,	US,	UZ,	VC,	VN	ı, YU,	ZA,	ZM,	ZW			
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			BF,	ВJ,	CF,	CG,	CI,	CM,	GA,	GN,	GΩ), GW,	ML,	MR,	NE,	SN,	TD,	TG
												2003-						
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		4113																
	EP	1553	411			A1		2005	0713		EP	2003-	7487	31		2	0031	006
		R:										, IT,						
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											JP	2004-	8365	4		2	0040	322
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PRIOR	RIT	APP:	LN.	INFO	. :							2002-						
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												2003-						
												2003-	JP12	793		W 2	0031	006
OTHER	2 50	TIRCE.	(8) .			MARI	PAT	140 •	3266,	41								

OTHER SOURCE(S): MARPAT 140:326641

AB Disclosed is an indicator for assessing body odor containing at least one member selected from the group consisting of (A) a

B-hydroxycarboxylic acid compound R1C(0H)(R2)CH2COOH (R1 = C1-4 alkyl; R2 = H, C1-4 alkyl; total number of C1 s < 10), (B) a derivative of the above compound A, (C) an alc. compound having a mercapto group at the 3-position HSC(R3) (R4)CH(R5)CH2OH (R3 = H, Me; R4 = C1-3 alkyl R5 = H, Meas; total number of C is \leq 8), and (D) a derivative of the above compound C, especially 3-hydroxy-3-methylhexanoic acid and 3-mercapto-3-methylhexanoi. Also it is intended to provide a process for producing the above compound C which comprises incubating human-origin perspiration in an environment with an oxygen concentration of 10 volume/volume8 or below. Further, it is

intended to provide a kit for assessing human body odor which contains a color developing reagent reacting with a human perspiration-origin B-hydroxycarboxylic acid.

IT 58888-76-9P, 3-Hydroxy-3-methylhexanoic acid

RL: ANT (Analyte); BSŪ (Biological study, unclassified); PUR (Purification or recovery); ANST (Analytical study); BIOL (Biological study); PREP (Preparation)

(indicator for assessing body odor containing β -hydroxy carboxylates or 3-mercapto alcs., process for producing same, body odor assessment method, method of assessing efficaciousness of deodorant and kit for conveniently assessing body odor)

RN 58888-76-9 CAPLUS

CN Hexanoic acid, 3-hvdroxv-3-methvl- (CA INDEX NAME)

OH n-Pr-C-CH2-CO2H Me

THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS REFERENCE COUNT: 3 RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L10 ANSWER 8 OF 8 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2003:122930 CAPLUS

DOCUMENT NUMBER: 139:2719

TITLE: A Specific Bacterial Aminoacylase Cleaves Odorant Precursors Secreted in the Human Axilla

Natsch, Andreas; Gfeller, Hans; Gygax, Peter; Schmid, AUTHOR(S):

Joachim; Acuna, Gonzalo Givaudan Duebendorf Ltd., Duebendorf, CH-8600, Switz. CORPORATE SOURCE:

SOURCE: Journal of Biological Chemistry (2003), 278(8),

5718-5727 CODEN: JBCHA3; ISSN: 0021-9258

PUBLISHER: American Society for Biochemistry and Molecular

Biology DOCUMENT TYPE: Journal LANGUAGE: English

Human axillary odor is known to be formed upon the action of

Corvnebacteria sp. on odorless axilla secretions. The known axilla odor determinant 3-methyl-2-hexenoic acid was identified in hydrolyzed axilla secretions along with a chemical related compound, 3-hydroxy-3-methylhexanoic acid. The natural precursors of both these acids were purified from non-hydrolyzed axilla secretions. From liquid chromatog./mass spectrometry anal., it appeared that the acids are covalently linked to a glutamine residue in fresh axilla secretions, and the corresponding conjugates were synthesized for confirmation. Bacterial isolates obtained from the human axilla and belonging to the Corynebacteria were found to release the acids from these odorless precursors in vitro. A Zn2+-dependent aminoacylase mediating this cleavage was purified from Corynebacterium striatum Ax20, and the corresponding gene agaA was cloned and heterologously expressed in Escherichia coli. The enzyme is highly specific for the glutamine residue but has a low specificity for the acvl part of the substrate. agaA is closely related to many genes coding for enzymes involved in the cleavage of N-terminal acyl and aryl substituents from amino acids. This is the first report of the structure elucidation of precursors for human body odorants and the isolation of the bacterial enzyme involved in their

cleavage. 58888-76-9, 3-Hvdroxv-3-methylhexanoic acid

RL: BSU (Biological study, unclassified); BIOL (Biological study) (anal. of hydrolyzed axilla secretions)

RN 58888-76-9 CAPLUS

CN Hexanoic acid, 3-hydroxy-3-methyl- (CA INDEX NAME)

REFERENCE COUNT:

THERE ARE 33 CITED REFERENCES AVAILABLE FOR THIS 33 RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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COST IN U.S. DOLLARS
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=> s 15 and 17 and 19 COMBINATION OF STRUCTURE AND TEXT TERMS NOT VALID The query entered contains both search terms created by structure-building or screen commands and text search terms. L#s created via the STRUCTURE or SCREEN commands must be searched in the structures files separately from text terms or profiles. The L# answer sets from structure searches can be used in crossover searches and can be combined with text terms.

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SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -6.56

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FILE COVERS 1907 - 24 Mar 2009 VOL 150 ISS 13 FILE LAST UPDATED: 23 Mar 2009 (20090323/ED)

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http://www.cas.org/legal/infopolicy.html

This file contains CAS Registry Numbers for easy and accurate substance identification.

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=> s 15 18977 L5 L13

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L14 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 2006:768974 CAPLUS
DOCUMENT NUMBER:
                         145:181018
TITLE:
                         Inhibition of sweat malodor
                         Starkenmann, Christian; Clark, Anthony; Troccaz,
INVENTOR(S):
                        Myriam; Niclass, Yvan
PATENT ASSIGNEE(S):
                        Firmenich SA, Switz.
SOURCE:
                         PCT Int. Appl., 36pp.
                         CODEN: PIXXD2
DOCUMENT TYPE:
                         Patent
LANGUAGE:
                         English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:
     PATENT NO.
                         KIND DATE APPLICATION NO. DATE
                         A2
     WO 2006079934
                         A2 20060803
A3 20061116
                                           WO 2006-IB50098
                                                                    20060111
     WO 2006079934
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             MZ, NA, NG, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE,
             SG, SK, SL, SM, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC,
             VN, YU, ZA, ZM, ZW
         RW: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE,
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             GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY,
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     EP 1846436
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                          A2 20071024
                                                                      20060111
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     CN 101107263
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                               20080116 CN 2006-80003050
                                                                     20060111
     US 20080025935
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                                20080131
                                           US 2007-779462
                                                                     20070718
     IN 2007KN02817
                          A
                               20070907
                                            IN 2007-KN2817
                                                                     20070801
PRIORITY APPLN. INFO.:
                                              EP 2005-100619
                                                                 A 20050131
                                              US 2005-647777P
                                                                 P 20050131
                                              WO 2006-IB50098
                                                                 W 20060111
OTHER SOURCE(S):
                         MARPAT 145:181018
    The present invention relates to a method for screening compds. having the
     ability to prevent, treat or reduce malodor development on body surfaces.
     In particular, the method allows to efficiently screen for compound having
     the ability of preventing sweat malodor development caused by volatile
     sulfur compds. (VSCs). The present invention is based on the finding of
     the direct precursor of naturally VSCs, which is present in human sweat
     and which will be metabolized by Staphylococci to VSCs.
REFERENCE COUNT:
                          3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS
                                RECORD, ALL CITATIONS AVAILABLE IN THE RE FORMAT
L14 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER:
                       2004:308604 CAPLUS
DOCUMENT NUMBER:
                          140:326641
TITLE:
                          Indicator for assessing body odor, process for
                          producing the same, body odor assessment method,
                         method of assessing efficaciousness of deodorant and
                         kit for conveniently assessing body odor
                         Yabuki, Masayuki; Hasegawa, Yoshihiro; Matsukane,
INVENTOR(S):
                         Masamoto; Yabe, Emi
PATENT ASSIGNEE(S):
                        Kao Corporation, Japan
                         PCT Int. Appl., 86 pp.
SOURCE:
                         CODEN: PIXXD2
DOCUMENT TYPE:
                         Patent
LANGUAGE:
                         Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:
    PATENT NO.
                        KIND DATE APPLICATION NO. DATE
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     WO 2004031766
                         A1 20040415 WO 2003-JP12793 20031006
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             GH, GM, HR, HU, ID, IL, IN, IS, KE, KG, KP, KR, KZ, LC, LK, LR,
             LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM,
             PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN,
             TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW
         RW: GH, GM, KE, LS, MM, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR,
             BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG

68776 Al 20040423 AU 2003-268776 20031006

109454 A 20041104 JP 2003-346596 20031006
     AU 2003268776
     JP 2004309454
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                         B2
                          B2 20080709
A1 20050713
     EP 1553411
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R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NI, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK JP 2005017272 A 20050120 JP 2004-83654 20040322

B2 20080423

JP 4081034

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US 20060160240 A1 20060720 US 2005-529897 20051216
PRIORITY APPLN. INFO:: JP 2002-293104 A 20021004
JP 2003-83801 A 20030325
JP 2003-116582 A 20030425
JP 2003-160882 A 20030604
JP 2003-346586 A 20031006
WO 2003-346586 A 20031006
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OTHER SOURCE(S): MARPAT 140:326641

AB Disclosed is an indicator for assessing body odor containing at least one

member selected from the group consisting of (A) a β-hydroxycarboxylic acid compound RIC(OH) (R2)CH2COOH (RI = Cl-4 alkyl; R2 = H, Cl-4 alkyl; total number of C is ≤ 10), (B) a derivative of the above compound A, (C) an alc. compound having a mercapto group at the 3-position HSC(R3) (R4)CH(R5)CH2OH (R3 = H, Me; R4 = Cl-3 alkyl R5 = H, Mesa; total number of C is ≤ 8), and (D) a derivative of the above compound C, especially 3-hydroxy-3-methylhexanoic acid and 3-mercapto-3-methylhexanol. Also it is intended to provide a process for producing the above compound C which comprises incubating human-origin perspiration in an environment with an oxygen concentration of 10 volume/volume8 or below. Further, it is

RECORD, ALL CITATIONS AVAILABLE IN THE RE FORMAT

intended to provide a kit for assessing human body odor which contains a color developing reagent reacting with a human perspiration-origin

 β -hydroxycarboxylic acid.

REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS

L14 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2002:594647 CAPLUS

DOCUMENT NUMBER: 137:145224

TITLE: An ATPase inhibitor-containing cosmetic products for

the reduction of sweat acidity

INVENTOR(S): Beck, Jonathan Samuel; Burry, Jason Shaun; Evans, Richard Livesey; Granger, Dominic; Laprade, Raynald;

Marsolais, Mireille

PATENT ASSIGNEE(S): Unilever PLC, UK; Unilever NV; Hindustan Lever Limited

SOURCE: PCT Int. Appl., 35 pp.

CODEN: PIXXD2
DOCUMENT TYPE: Patent

LANGUAGE: English FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT				KIN	D	DATE			APPL	ICAT	ION :	NO.		D.	ATE	
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WO 200	20604	02		A1		2002	0808		WO 2	002-	EP67	0		2	0020	121
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AU 200	22280	60		A1		2002	0812		AU 2	002-	2280	60		2	0020	121
US 200	20146	376		A1		2002	1010		US 2	002-	6618	3		2	0020	131
US 650	9010			B2		2003	0121									
PRIORITY AP	PLN.	INFO	. :						GB 2	001-	2562			A 2	0010	201
									WO 2	002-	EP67	0	1	ii 2	0020	121
OTHER SOURC	E(S):			MAR	PAT	137:	1452	24								

AB A cosmetic method of reducing the acidity of sweat excreted from human eccrine glands comprises the topical application of a vacuolar-type

H+-ATPase (V-ATPase) inhibitor to the skin in the vicinity of the eccrine glands. The method may result in a range of benefits, including enhanced appreciation of topically-applied perfume and enhanced efficacy of topically-applied antiperspirant salt. Cosmetic products and compnss. comprising a V-ATPase inhibitor and selected other components are also claimed. For example, olygomycin (at 20 µg/mL), bafilomycin A1 (at 6.2 µg/mL), and concanamycin A (at 0.1 µg/mL) all inhibit proton transfer out of the cells of the reabsorptive duct affecting the pH recovery by 12, 27, and 5%, resp., compared to 100% pH recovery in the control (no V-ATPase inhibitor).

REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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L15 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2009 ACS on STN
    2004:308604 CAPLUS
DN
    140:326641
ED
    Entered STN: 15 Apr 2004
ΤТ
     Indicator for assessing body odor, process for producing the same, body
     odor assessment method, method of assessing efficaciousness of deodorant
     and kit for conveniently assessing body odor
    Yabuki, Masayuki; Hasegawa, Yoshihiro; Matsukane, Masamoto; Yabe, Emi
IN
PA
    Kao Corporation, Japan
     PCT Int. Appl., 86 pp.
SO
     CODEN: PIXXD2
DT
     Patent
LA
    Japanese
     ICM G01N033-497
IC
     ICS G01N033-15; C11B009-00; A61K007-46
     62-4 (Essential Oils and Cosmetics)
     Section cross-reference(s): 9
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CLASS
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US 20060160240 IPCI
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                       A61K008/02; A61Q015/00; C09B026/02; G01N033/497; K61K
OS
    MARPAT 140:326641
AB
    Disclosed is an indicator for assessing body odor containing at least one
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member selected from the group consisting of (A) as β -hydroxycarboxylic acid compound RIC(OH)(R2)CH2COOH (R1 = C1-4 alkyl; total number of C is \leq 10), (B) a derivative of the

above compound A, (C) an alc. compound having a mercapto group at the 3-position HSC(R3)(R4)CH(R5)CH2OH (R3 = H, Me; R4 = C1-3 alkyl R5 = H, Meas; total number of C is \leq 8), and (D) a derivative of the above compound C, especially 3-hydroxy-3-methylhexanoic acid and 3-mercapto-3-methylhexanol. Also it is intended to provide a process for producing the above compound C which comprises incubating human-origin perspiration in an environment with an oxygen concentration of 10 volume/volume% or below. Further, it is

to provide a kit for assessing human body odor which contains a color developing reagent reacting with a human perspiration-origin

B-hydroxycarboxylic acid. ST hydroxy carboxylate body odor indicator; mercapto alc body odor indicator;

deodorant assessment hydroxy carboxylate mercapto alc

тт Carboxylic acids, biological studies

RL: ANT (Analyte); BSU (Biological study, unclassified); PUR (Purification or recovery); ANST (Analytical study); BIOL (Biological study); PREP (Preparation)

(hydroxy; indicator for assessing body odor containing β-hydroxy carboxylates or 3-mercapto alcs., process for producing same, body odor assessment method, method of assessing efficaciousness of deodorant and kit for conveniently assessing body odor)

Deodorants (personal)

Human

Odor and Odorous substances

Sweat.

(indicator for assessing body odor containing \(\beta\)-hydroxy carboxylates or 3-mercapto alcs., process for producing same, body odor assessment method, method of assessing efficaciousness of deodorant and kit for conveniently assessing body odor)

Alcohols, biological studies

RL: ANT (Analyte); BSU (Biological study, unclassified); PUR (Purification or recovery); ANST (Analytical study); BIOL (Biological study); PREP (Preparation)

(indicator for assessing body odor containing \(\beta\)-hydroxy carboxylates or 3-mercapto alcs., process for producing same, body odor assessment method, method of assessing efficaciousness of deodorant and kit for conveniently assessing body odor)

51755-83-0P, 3-Mercaptohexanol 58888-76-9P, 3-Hydroxy-3-methylhexanoic 227456-27-1P 227456-33-9P 307964-23-4P 548740-99-4P RL: ANT (Analyte); BSU (Biological study, unclassified); PUR (Purification or recovery); ANST (Analytical study); BIOL (Biological study); PREP (Preparation)

(indicator for assessing body odor containing 6-hydroxy carboxylates or 3-mercapto alcs., process for producing same, body odor assessment method, method of assessing efficaciousness of deodorant and kit for conveniently assessing body odor)

10401-59-9, 9-Anthryldiazomethane

RL: ARG (Analytical reagent use); ANST (Analytical study); USES (Uses) (indicator for assessing body odor containing \(\beta\)-hydroxy carboxylates or 3-mercapto alcs., process for producing same, body odor assessment method, method of assessing efficaciousness of deodorant and kit for conveniently assessing body odor)

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